

# High Resolution FTIR Spectrum of Chlorofluoroethyne, FCCl, below 1000 cm<sup>-1</sup>. Analysis of the $\nu_3$ , $\nu_4$ , $\nu_5$ , $2\nu_4$ , $\nu_4 + \nu_5$ and $2\nu_5$ Bands, and *ab initio* Calculations

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High resolution infrared spectra of FCCl have been measured and analyzed by polynomial methods. In the region below 350 cm<sup>-1</sup>, the analysis is straightforward and yields parameters for the  $\nu_4 = 1$  and  $\nu_5 = 1$  states. Between 350 and 800 cm<sup>-1</sup> there are strong anharmonic interactions in the  $2\nu_5/\nu_4 + \nu_5/\nu_3/2\nu_4$  tetrad which have been unravelled with the use of a model that employs *ab initio* interaction constants. Observed and theoretically predicted wavenumbers are in excellent agreement for all bands studied.

*Key words:* Infrared Spectrum; High Resolution; Alkyne; Fermi Resonance; *ab initio* Calculations.

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